

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092120\
 Data File : VN063514.D
 Acq On : 22 Sep 2020 3:42
 Operator : JC/MD
 Sample : L4082-09
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 RE134D2-20200915

Manual Integrations
 APPROVED

MMDadoda
 9/22/2020 2:08:27 PM

Quant Time: Sep 22 05:57:16 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 16 13:05:20 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	228176	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	425325	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	428292	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	180142	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	189991	48.91	ug/l	0.00
Spiked Amount	50.000		Recovery	=	97.82%	
35) Dibromofluoromethane	7.55	113	139571	47.40	ug/l	0.00
Spiked Amount	50.000		Recovery	=	94.80%	
50) Toluene-d8	10.06	98	544110	49.14	ug/l	0.00
Spiked Amount	50.000		Recovery	=	98.28%	
62) 4-Bromofluorobenzene	12.38	95	215337	53.06	ug/l	0.00
Spiked Amount	50.000		Recovery	=	106.12%	
Target Compounds						
3) Chloromethane	2.04	50	2169	0.634	ug/l	97
9) 1,1,2-Trichlorotrifluoroet	3.72	101	61090	24.030	ug/l	99
27) cis-1,2-Dichloroethene	6.79	96	7708m	2.387	ug/l	
44) Trichloroethene	8.80	130	406899	129.696	ug/l	98
49) 1,4-Dioxane	9.17	88	2327	41.504	ug/l #	82
64) Tetrachloroethene	10.60	164	45122	15.472	ug/l	95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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